

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110822\
 Data File : BF131072.D
 Acq On : 08 Nov 2022 17:58
 Operator : CG\JU
 Sample : N5415-11RE
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 R-24RE

Quant Time: Nov 09 04:07:31 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 08 01:40:43 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.898	152	78351	20.000	ng	0.00
21) Naphthalene-d8	8.186	136	272468	20.000	ng	# 0.00
39) Acenaphthene-d10	9.945	164	126446	20.000	ng	0.00
64) Phenanthrene-d10	11.433	188	214220	20.000	ng	0.00
76) Chrysene-d12	14.086	240	228755	20.000	ng	0.00
86) Perylene-d12	15.592	264	222644	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	377607	77.713	ng	0.00
7) Phenol-d6	6.522	99	440536	74.281	ng	0.00
23) Nitrobenzene-d5	7.463	82	288812	49.133	ng	0.00
42) 2,4,6-Tribromophenol	10.733	330	93055	66.843	ng	0.00
45) 2-Fluorobiphenyl	9.257	172	453066	48.831	ng	-0.01
79) Terphenyl-d14	13.027	244	478360	33.526	ng	0.00
Target Compounds						
87) Indeno(1,2,3-cd)pyrene	17.098	276	10540	3.847	ng	Qvalue 95
92) Benzo(g,h,i)perylene	17.568	276	10926	4.042	ng	# 95

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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